=> d scan

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2-(ethylamino)-2,3-dihydro-, (1R,2R)-rel- (9CI)

MF C11 H15 N O

CI COM

Relative stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):50

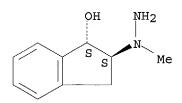
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)- (9CI)

MF C10 H14 N2 O

CI COM

Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-,

(2Z)-2-butenedioate (1:1) (salt) (9CI)

MF C10 H14 N2 O . C4 H4 O4

CM 1

Absolute stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN Butanedioic acid, compd. with (1R,2R)-2,3-dihydro-2-(1-methylhydrazino)-1H-inden-1-ol (1:2) (9CI)

MF C10 H14 N2 O . 1/2 C4 H6 O4

CM 1

Absolute stereochemistry.

CM 2

 $HO_2C-CH_2-CH_2-CO_2H$

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

MF C10 H14 N2 O . 1/2 C4 H6 O6

CM 1

Absolute stereochemistry.

CM 2

Absolute stereochemistry.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

MF C11 H16 N2 O . C4 H4 O4

CM 1

Relative stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2-(ethylnitrosoamino)-2,3-dihydro-, (1R,2R)-rel- (9CI)

MF C11 H14 N2 O2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylnitrosoamino)-, (1R,2R)-rel- (9CI)
MF C10 H12 N2 O2

Relative stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2S)-,
(2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C10 H14 N2 O . C4 H4 O4

CM 1

Absolute stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 2-Butenedioic acid (2E) - (9CI)

MF C4 H4 O4

CI COM

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-6-methoxy-2-(methylamino)-, (1R,2R)-rel- (9CI)

MF C11 H15 N O2

CI COM

Relative stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylamino)-, (1S-trans)- (9CI)

MF C10 H13 N O

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-rel-,

(2Z)-2-butenedioate (1:1) (salt) (9CI) MF C10 H14 N2 O . C4 H4 O4

CM :

Relative stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN Butanedioic acid, compd. with (1S,2S)-2,3-dihydro-2-(1-methylhydrazino)-1Hinden-1-ol (1:2) (9CI)

MF C10 H14 N2 O . 1/2 C4 H6 O4

CM 1

Absolute stereochemistry. Rotation (+).

CM 2

 ${\tt HO_2C-CH_2-CH_2-CO_2H}$

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)-, (2R,3R)-2,3-dihydroxybutanedioate (2:1) (salt) (9CI)
MF C10 H14 N2 O . 1/2 C4 H6 O6

CM 1

Absolute stereochemistry. Rotation (+).

CM 2

Absolute stereochemistry.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2S)-rel-,

(2Z)-2-butenedioate (1:1) (salt) (9CI)

MF C10 H14 N2 O . C4 H4 O4

CM 1

Relative stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylnitrosoamino)-, (1S,2S)- (9CI)

MF C10 H12 N2 O2

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-5-methyl-2-(methylnitrosoamino)-, (1R,2R)-rel(9CI)

MF C11 H14 N2 O2

Relative stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-6-methyl-2-(methylamino)-, (1R,2R)-rel- (9CI)
MF C11 H15 N O

Relative stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 2-Butenedioic acid (2Z) - (9CI)

MF C4 H4 O4

CI COM

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN Oxidase, monoamine (9CI)

MF Unspecified

CI MAN

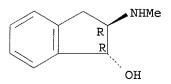
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylamino)-, (1R-trans)- (9CI)

MF C10 H13 N O

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2-(1-ethylhydrazino)-2,3-dihydro-, (1R,2R)-rel-,

(2Z)-2-butenedioate (1:1) (salt) (9CI)

MF C11 H16 N2 O . C4 H4 O4

CM 1

CM

Double bond geometry as shown.

REGISTRY COPYRIGHT 2003 ACS on STN L7

1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-, IN (2E) -2-butenedioate (2:1) (salt) (9CI)

ΜF C10 H14 N2 O . 1/2 C4 H4 O4

CM

Absolute stereochemistry.

CM 2

Double bond geometry as shown.

41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN L7

1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-, IN (2S, 3S) -2, 3-dihydroxybutanedioate (2:1) (salt) (9CI)

C10 H14 N2 O . 1/2 C4 H6 O6

CM1

MF

Absolute stereochemistry.

2 CM

Absolute stereochemistry.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2S)- (9CI)

MF C10 H14 N2 O

CI COM

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

MF C11 H16 N2 O2 . C4 H4 O4

CM 1

Relative stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylnitrosoamino)-, (1R,2S)- (9CI)

MF C10 H12 N2 O2

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-5-methyl-2-(methylamino)-, (1R,2R)-rel- (9CI)

MF C11 H15 N O

Relative stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN Butanedioic acid (9CI)

MF C4 H6 O4

CI COM

 ${\tt HO_2C-CH_2-CH_2-CO_2H}$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1-Oxa-2-azaspiro[2.5]octane (7CI, 8CI, 9CI)

MF C6 H11 N O

CI RPS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN VAP 1 (9CI)

MF Unspecified

CI MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

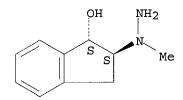
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)-,

(2Z)-2-butenedioate (1:1) (salt) (9CI)

MF C10 H14 N2 O . C4 H4 O4

CM 1

Absolute stereochemistry. Rotation (+).



CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

MF C10 H14 N2 O . 1/2 C4 H4 O4

CM 1

Absolute stereochemistry. Rotation (+).

CM 2

Double bond geometry as shown.

CM 1

Absolute stereochemistry. Rotation (+).

CM 2

Absolute stereochemistry.

Absolute stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-6-methyl-2-(1-methylhydrazino)-, (1R,2R)-rel-,

(2Z)-2-butenedioate (1:1) (salt) (9CI)

MF C11 H16 N2 O . C4 H4 O4

CM 1

Relative stereochemistry.

CM 2

Double bond geometry as shown.

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-2-amine, 2,3-dihydro-1-methoxy-N-methyl-N-nitroso-, (1S,2S)-

(9CI)

MF C11 H14 N2 O2

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylamino)-, (1R,2S)- (9CI)

MF C10 H13 N O

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN Butanedioic acid, 2,3-dihydroxy- (2R,3R)- (9CI)

MF C4 H6 O6

CI COM

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

IN Butanedioic acid, 2,3-dihydroxy-, (2S,3S)- (9CI)

MF C4 H6 O6

CI COM

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=>

=> d bib hitstr 2-7

L14 ANSWER 2 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN

AN 1996:332010 CAPLUS

DN 125:86465

TI Synthesis and stereochemistry of oxazolo[4',5':1,2]benz[5,4,3-c,d]indole derivatives as intermediates on the way to a selective dopamine autoreceptor agonist

AU Gmeiner, Peter; Bollinger, Bernd; Lotter, Hermann

CS Pharmazeutisches Inst. der Univ. Bonn, Bonn, D-53121, Germany

SO Journal of Heterocyclic Chemistry (1996), 33(2), 481-483 CODEN: JHTCAD; ISSN: 0022-152X

PB HeteroCorporation

DT Journal

LA English

IT 178551-99-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(prepn. of (dipropylamino)benz[cd]indolone)

RN 178551-99-0 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-[1,2,2a,3,4,5-hexahydro-5-hydroxy-2-oxobenz[cd]indol-4-yl]-, bis(phenylmethyl) ester, (2a.alpha.,4.beta.,5.beta.)- (9CI) (CA INDEX NAME)

Relative stereochemistry.

L14 ANSWER 3 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN

AN 1995:801122 CAPLUS

DN 124:8665

TI Synthesis, pharmacological investigation and computational studies on a tricyclic ergoline analog with selective dopamine autoreceptor activity

AU Gmeiner, Peter; Bollinger, Bernd; Mierau, Joachim; Hoefner, Georg

CS Pharmazeutisches Institut, Universitaet Bonn, Bonn, D-53121, Germany

SO Archiv der Pharmazie (Weinheim, Germany) (1995), 328(7-8), 609-14 CODEN: ARPMAS; ISSN: 0365-6233

PB VCH

DT Journal

LA English

OS CASREACT 124:8665

IT 171006-44-3P

RL: SPN (Synthetic preparation); PREP (Preparation) (synthesis, pharmacol. investigation and computational studies on a tricyclic ergoline analog with selective dopamine autoreceptor activity)

RN 171006-44-3 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-[1,2a-bis(diethoxymethyl)-1,2,2a,3,4,5-hexahydro-5-hydroxy-2-oxobenz[cd]indol-4-yl]-, bis(phenylmethyl) ester, (2a.alpha.,4.beta.,5.alpha.)- (9CI) (CA INDEX NAME)

Relative stereochemistry.

L14 ANSWER 4 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN

AN 1992:426035 CAPLUS

DN 117:26035

TI Synthesis of amines and amino alcohols by electrophilic amination and highly stereoselective reduction

AU Gmeiner, Peter; Bollinger, Bernd

CS Inst. Pharm. Lebensmittelchem., Univ. Muenchen, Munich, 8000/2, Germany

SO Liebigs Annalen der Chemie (1992), (3), 273-8 CODEN: LACHDL; ISSN: 0170-2041

DT Journal

LA English

OS CASREACT 117:26035

IT 138206-94-7P 138206-95-8P 138408-15-8P

138408-16-9P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(prepn. and hydrogenolysis of)

RN 138206-94-7 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-2-naphthalenyl)-, bis(phenylmethyl) ester, cis-(9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 138206-95-8 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-7-methoxy-

2-naphthalenyl)-, bis(phenylmethyl) ester, cis- (9CI) (CA INDEX NAME) Relative stereochemistry.

RN 138408-15-8 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-2-naphthalenyl)-, bis(phenylmethyl) ester, trans- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 138408-16-9 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-7-methoxy-2-naphthalenyl)-, bis(phenylmethyl) ester, trans- (9CI) (CA INDEX NAME)

Relative stereochemistry.

IT 138408-13-6P 138408-14-7P 138408-17-0P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(prepn. and nitrogen-nitrogen bond cleavage of)

RN 138408-13-6 CAPLUS

CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-, monohydrochloride, cis-(9CI) (CA INDEX NAME)

● HCl

RN 138408-14-7 CAPLUS

CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-7-methoxy-, monohydrochloride, cis- (9CI) (CA INDEX NAME)

Relative stereochemistry.

● HCl

RN 138408-17-0 CAPLUS

CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-, monohydrochloride, trans-(9CI) (CA INDEX NAME)

Relative stereochemistry.

● HCl

RN 138408-18-1 CAPLUS

CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-7-methoxy-, monohydrochloride, trans- (9CI) (CA INDEX NAME)

● HCl

L14 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN

AN 1992:41818 CAPLUS

DN 116:41818

TI Efficient methodology for the preparation of .beta.-aminotetralin derivatives via electrophilic amination

AU Gmeiner, Peter; Bollinger, Bernd

CS Inst. Pharm. Lebensmittelchem., Ludwig-Maximilians-Univ., Munich, 8000/2, Germany

SO Tetrahedron Letters (1991), 32(42), 5927-30 CODEN: TELEAY; ISSN: 0040-4039

DT Journal

LA English

OS CASREACT 116:41818

IT 138206-94-7P 138206-95-8P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(prepn. and transesterification-cyclization of)

RN 138206-94-7 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-2-naphthalenyl)-, bis(phenylmethyl) ester, cis- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 138206-95-8 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-7-methoxy-2-naphthalenyl)-, bis(phenylmethyl) ester, cis-(9CI) (CA INDEX NAME)

L14 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN

AN 1970:121454 CAPLUS

DN 72:121454

TI Photochemical and thermal 1,2- and 1,4-cycloaddition reactions of azodicarbonyl compounds with monoolefins

AU Koerner von Gustorf, Ernst; White, Danny V.; Kim, Bongsub; Hess, Dieter; Leitich, Johannes

CS Abt. Strahlenchem., Max Planck Inst. Kohlenforsch., Muelheim, Fed. Rep. Ger.

SO Journal of Organic Chemistry (1970), 35(4), 1155-65 CODEN: JOCEAH; ISSN: 0022-3263

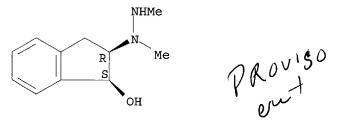
DT Journal

LA English

RN 23358-19-2 CAPLUS

CN 1-Indanol, 2-(1,2-dimethylhydrazino)-, cis- (8CI) (CA INDEX NAME)

Relative stereochemistry.



RN 23358-23-8 CAPLUS

CN Bicarbamic acid, (1-hydroxy-2-indanyl)-, diethyl ester (8CI) (CA INDEX NAME)

L14 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN AN 1970:121451 CAPLUS

DN 72:121451

TI Azodiformate adduct of indene and the stereochemistry of some 1,2-disubstituted indans

AU Huebner, Charles F.; Donoghue, Ellen M.; Novak, Carol J.; Dorfman, Louis; Wenkert, Ernest

CS Chem. Res. Div., CIBA Pharm. Co., Summit, NJ, USA

SO Journal of Organic Chemistry (1970), 35(4), 1149-54 CODEN: JOCEAH; ISSN: 0022-3263

DT Journal

LA English

IT 23358-19-2P 23359-96-8P

RN 23358-19-2 CAPLUS

CN 1-Indanol, 2-(1,2-dimethylhydrazino)-, cis- (8CI) (CA INDEX NAME)

Relative stereochemistry.

RN 23359-96-8 CAPLUS

=>

CN Acetic acid, 2-(1-hydroxy-2-indanyl)-1,2-dimethylhydrazide acetate (ester) (8CI) (CA INDEX NAME)